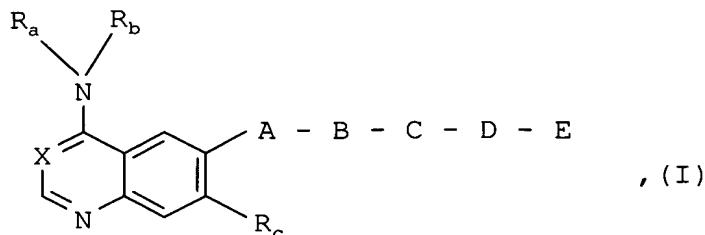


Amendments to the Claims:

This listing of claims will replace all prior versions and listing of claims in the application:

Listing of Claims:

1. (currently amended) A quinazoline compound of formula



wherein

R_a denotes a hydrogen atom or a C_{1-4} -alkyl group,

R_b denotes a phenyl, benzyl or 1-phenylethyl group wherein the phenyl nucleus is substituted in each case by the groups R_1 to R_3 , whilst

R_1 and R_2 , which may be identical or different, in each case denote a hydrogen, fluorine, chlorine, bromine or iodine atom,

a C_{1-4} -alkyl, hydroxy, C_{1-4} -alkoxy, C_{3-6} -cycloalkyl, C_{4-6} -cycloalkoxy, C_{2-5} -alkenyl or C_{2-5} -alkynyl group,

an aryl, aryloxy, arylmethyl or arylmethoxy group,

a C_{3-5} -alkenyloxy or C_{3-5} -alkynyoxy group, whilst the unsaturated moiety may not be linked to the oxygen atom,

a C_{1-4} -alkylsulfenyl, C_{1-4} -alkylsulfinyl, C_{1-4} -alkylsulfonyl, C_{1-4} -alkylsulfonyloxy, trifluoromethylsulfenyl, trifluoromethylsulfinyl or trifluoromethylsulfonyl group,

a methyl or methoxy group substituted by 1 to 3 fluorine atoms,

an ethyl or ethoxy group substituted by 1 to 5 fluorine atoms,

a cyano or nitro group or an amino group optionally substituted by one or two C₁₋₄-alkyl groups, wherein the substituents may be identical or different, or

R₁ together with R₂, if they are bound to adjacent carbon atoms, denote a -CH=CH-CH=CH group and

R₃ denotes a hydrogen, fluorine, chlorine or bromine atom,

a C₁₋₄-alkyl, trifluoromethyl or C₁₋₄-alkoxy group,

X denotes a nitrogen atom,

A denotes an imino group optionally substituted by a C₁₋₄-alkyl group,

B denotes a carbonyl group,

C denotes a ~~1,3-allenylene~~ -CH=C=CH-, ~~1,1->C=CH₂~~ or ~~1,2-vinylene~~ -CH=CH- group which may be substituted in each case by one or two methyl groups or by a trifluoromethyl group,

an ~~ethynylene~~ -C≡C- group or

a ~~1,3-butadien-1,4-ylene~~ -CH=CH-CH=CH- group optionally substituted by 1 to 4 methyl groups or by a trifluoromethyl group,

D denotes an alkylene group wherein the alkylene moiety contains 1 to 8 carbon atoms and additionally 1 to 4 hydrogen atoms in the alkylene moiety may be replaced by fluorine atoms,

E denotes an amino, C₁₋₄-alkylamino or di-(C₁₋₄-alkyl)-amino group wherein the alkyl moieties may be identical or different,

a C_{2-4} -alkylamino group wherein the alkyl moiety is substituted in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , whilst

R_5 denotes a hydroxy, C_{1-4} -alkoxy, amino, C_{1-4} -alkylamino or di- $(C_{1-4}$ -alkyl)-amino group,

an $N-(C_{1-4}$ -alkyl)- $N-(C_{2-4}$ -alkyl)-amino group wherein the C_{2-4} -alkyl moiety is substituted in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , whilst R_5 is as hereinbefore defined,

a di- $(C_{2-4}$ -alkyl)-amino group wherein the two C_{2-4} -alkyl moieties are substituted in each case in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , whilst the substituents may be identical or different and R_5 is as hereinbefore defined,

a C_{3-7} -cycloalkylamino or C_{3-7} -cycloalkyl- C_{1-3} -alkylamino group wherein in each case the nitrogen atom may be substituted by a further C_{1-4} -alkyl group,

R_c denotes a C_{4-7} -cycloalkoxy or C_{3-7} -cycloalkyl- C_{1-6} -alkoxy group wherein the cycloalkyl moiety in each case may be substituted by a C_{1-3} -alkyl, hydroxy, C_{1-4} -alkoxy, amino, C_{1-4} -alkylamino, di- $(C_{1-4}$ -alkyl)-amino, hydroxy- C_{1-2} -alkyl, C_{1-4} -alkoxy- C_{1-2} -alkyl, amino- C_{1-2} -alkyl, C_{1-4} -alkylamino- C_{1-2} -alkyl, or di- $(C_{1-4}$ -alkyl)-amino- C_{1-2} -alkyl group, whilst the abovementioned monosubstituted cycloalkyl moieties may additionally be substituted by a C_{1-3} -alkyl group,

whilst

by the aryl moieties mentioned in the definition of the abovementioned groups is meant a phenyl group which in each case may be monosubstituted by R_7 , mono-, di- or trisubstituted by R_8 or monosubstituted by R_7 and additionally mono- or disubstituted by R_8 , wherein the substituents may be identical or different and

R_7 denotes a cyano, carboxy, C_{1-4} -alkoxycarbonyl, aminocarbonyl, C_{1-4} -alkylaminocarbonyl, di- $(C_{1-4}$ -alkyl)-aminocarbonyl, C_{1-4} -alkylsulfonyl, C_{1-4} -alkylsulfinyl, C_{1-4} -alkylsulfonyl, hydroxy, C_{1-4} -alkylsulfonyloxy, trifluoromethyloxy, nitro, amino, C_{1-4} -alkylamino, di- $(C_{1-4}$ -alkyl)-amino, C_{1-4} -alkylcarbonylamino, N -(C_{1-4} -alkyl)- C_{1-4} -alkylcarbonylamino, C_{1-4} -alkylsulfonylamino, N -(C_{1-4} -alkyl)- C_{1-4} -alkylsulfonylamino, aminosulfonyl, C_{1-4} -alkylaminosulfonyl or di- $(C_{1-4}$ -alkyl)-aminosulfonyl group, and

R_8 denotes a fluorine, chlorine, bromine or iodine atom, a C_{1-4} -alkyl, trifluoromethyl or C_{1-4} -alkoxy group or

two groups R_8 , if they are bound to adjacent carbon atoms, together denote a C_{3-5} -alkylene or 1,3-butadien-1,4-ylene group,

or the tautomers, or stereoisomers or pharmaceutically acceptable salts thereof.

2. (currently amended) A quinazoline of formula I according to claim 1, wherein

R_a denotes a hydrogen atom,

R_b denotes a phenyl, benzyl or 1-phenylethyl group wherein the phenyl nucleus is substituted in each case by the groups R_1 to R_3 , whilst

R_1 and R_2 , which may be identical or different, in each case denote a hydrogen, fluorine, chlorine, bromine or iodine atom,

a C_{1-4} -alkyl, hydroxy, C_{1-4} -alkoxy, C_{3-6} -cycloalkyl, C_{4-6} -cycloalkoxy, C_{2-5} -alkenyl or C_{2-5} -alkynyl group,

an aryl, aryloxy, arylmethyl or arylmethoxy group,

a methyl or methoxy group substituted by 1 to 3 fluorine atoms,

a cyano or nitro group and

R_3 denotes a hydrogen, fluorine, chlorine or bromine atom,

a C_{1-4} -alkyl, trifluoromethyl or C_{1-4} -alkoxy group,

X denotes a nitrogen atom,

A denotes an imino group,

B denotes a carbonyl group,

C denotes a ~~1,3-allenylene~~ $-CH=C=CH-$, ~~1,1->C=CH₂~~ or ~~1,2-vinylene~~ $-CH=CH-$ group,

an ~~ethynylene~~ $-C\equiv C-$ or ~~1,3-butadien-1,4-ylene~~ $-CH=CH-CH=CH-$ group,

D denotes an alkylene group wherein the alkylene moiety in each case contains 1 to 4 carbon atoms and additionally 1 to 4 hydrogen atoms in the alkylene moiety may be replaced by fluorine atoms,

E denotes a di- $(C_{1-4}$ -alkyl)-amino group wherein the alkyl moieties may be identical or different,

an $N-(C_{1-4}$ -alkyl)- $N-(C_{2-4}$ -alkyl)-amino group wherein the C_{2-4} -alkyl moiety is substituted in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , where

R_5 denotes a hydroxy, C_{1-4} -alkoxy or di- $(C_{1-4}$ -alkyl)-amino group,

a di- $(C_{2-4}$ -alkyl)-amino group wherein the two C_{2-4} -alkyl moieties in each case are substituted in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , wherein the substituents may be identical or different and R_5 is as hereinbefore defined,

a C₃₋₇-cycloalkylamino or C₃₋₇-cycloalkyl-C₁₋₃-alkylamino group wherein in each case the nitrogen atom is substituted by a further C₁₋₄-alkyl group,

R_c denotes a C₄₋₇-cycloalkoxy or C₃₋₇-cycloalkyl-C₁₋₆-alkoxy group wherein the cycloalkyl moiety in each case may be substituted by a C₁₋₃-alkyl, hydroxy, C₁₋₄-alkoxy, di-(C₁₋₄-alkyl)-amino, hydroxy-C₁₋₂-alkyl, C₁₋₄-alkoxy-C₁₋₂-alkyl, or di-(C₁₋₄-alkyl)-amino-C₁₋₂-alkyl group, whilst the abovementioned monosubstituted cycloalkyl moieties may additionally be substituted by a C₁₋₃-alkyl group, , whilst

by the aryl moieties mentioned in the definition of the abovementioned groups is meant a phenyl group which may in each case be monosubstituted by R₇, mono-, di- or trisubstituted by R₈ or monosubstituted by R₇ and additionally mono- or disubstituted by R₈, wherein the substituents may be identical or different and

R₇ denotes a cyano, carboxy, C₁₋₄-alkoxycarbonyl, aminocarbonyl, C₁₋₄-alkylaminocarbonyl, di-(C₁₋₄-alkyl)-aminocarbonyl, C₁₋₄-alkylsulfonyl, C₁₋₄-alkylsulfinyl, C₁₋₄-alkylsulfonyl, hydroxy, C₁₋₄-alkylsulfonyloxy, trifluoromethoxy, nitro, amino, C₁₋₄-alkylamino, di-(C₁₋₄-alkyl)-amino, C₁₋₄-alkylcarbonylamino, N-(C₁₋₄-alkyl)-C₁₋₄-alkylcarbonylamino, C₁₋₄-alkylsulfonylamino, N-(C₁₋₄-alkyl)-C₁₋₄-alkylsulfonylamino, aminosulfonyl, C₁₋₄-alkylaminosulfonyl or di-(C₁₋₄-alkyl)-aminosulfonyl group, and

R₈ denotes a fluorine, chlorine, bromine or iodine atom, a C₁₋₄-alkyl, trifluoromethyl or C₁₋₄-alkoxy group or

two groups R₈, if they are bound to adjacent carbon atoms, together denote a C₃₋₅-alkylene or 1,3-butadien-1,4-ylene group,

or the tautomers, or stereoisomers or pharmaceutically acceptable salts thereof.

3. (currently amended) A quinazoline of formula I according to claim 1, wherein

R_a denotes a hydrogen atom,

R_b denotes a phenyl, benzyl or 1-phenylethyl group wherein the phenyl nucleus is substituted in each case by the groups R_1 and R_2 , where

R_1 and R_2 , which may be identical or different, in each case denote a hydrogen, fluorine, chlorine or bromine atom,

a methyl, trifluoromethyl or methoxy group,

X denotes a nitrogen atom,

A denotes an imino group,

B denotes a carbonyl group,

C denotes a ~~1,2-vinylene~~- $\text{CH}=\text{CH-}$ group,

an ~~ethynylene~~ $\text{-C}\equiv\text{C-}$ or ~~1,3-butadien-1,4-ylene~~ $\text{-CH}=\text{CH-CH}=\text{CH-}$ group,

D denotes a C_{1-4} -alkylene group,

E denotes a di- $(\text{C}_{1-4}\text{-alkyl})$ -amino group wherein the alkyl moieties may be identical or different,

an $\text{N-}(\text{C}_{1-4}\text{-alkyl})\text{-N-}(\text{C}_{2-4}\text{-alkyl})$ -amino group wherein the C_{2-4} -alkyl moiety is substituted in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , whilst

R_5 denotes a hydroxy, C_{1-3} -alkoxy or di- $(\text{C}_{1-3}\text{-alkyl})$ -amino group,

a di- $(\text{C}_{2-4}\text{-alkyl})$ -amino group wherein the two C_{2-4} -alkyl moieties in each case are substituted in β -, γ -, or δ -position with regard to the nitrogen atom of the amino group by the group R_5 , wherein the substituents may be identical or different and R_5 is as hereinbefore defined,

a C₃₋₅-cycloalkylamino or C₃₋₅-cycloalkyl-C₁₋₃-alkylamino group wherein in each case the nitrogen atom is substituted by a further C₁₋₃-alkyl group,

R_c denotes a C₄₋₇-cycloalkoxy or C₃₋₇-cycloalkyl-C₁₋₄-alkoxy group wherein the cycloalkyl moiety in each case may be substituted by a C₁₋₃-alkyl or C₁₋₃-alkoxy group, or the tautomers, or stereoisomers or pharmaceutically acceptable salts thereof.

4. (currently amended) A quinazoline of formula I according to claim 1, wherein

R_a denotes a hydrogen atom,

R_b denotes a phenyl, benzyl or 1-phenylethyl group, whilst the phenyl nucleus is substituted in each case by the radicals R₁ and R₂, whilst

R₁ and R₂, which may be identical or different, each denotes a hydrogen, fluorine, chlorine or bromine atom,

X denotes a nitrogen atom,

A denotes an imino group,

B denotes a carbonyl group,

C denotes a 1,2-vinylene -CH=CH-, ethynylene -C≡C- or 1,3-butadien-1,4-ylene -CH=CH-
CH=CH- group, |

D denotes an C₁₋₃-alkylene group,

E denotes a Di-(C₁₋₄-alkyl)-amino group, wherein the alkyl moieties may be identical or different,

a methylamino or ethylamino group each substituted at the nitrogen atom by a 2-methoxyethyl, 1-methoxy-2-propyl, 2-methoxypropyl, 3-methoxypropyl, cyclopropyl or cyclopropylmethyl group,

a bis-(2-methoxyethyl)amino group,

R_c denotes a cyclopropylmethoxy, cyclobutylmethoxy, cyclopentylmethoxy or cyclohexylmethoxy group,

a cyclobutoxy, cyclopentyloxy or cyclohexyloxy group,

or the tautomers, or stereoisomers or pharmaceutically acceptable salts thereof.

5. (currently amended) A quinazoline of formula I according to claim 1, wherein

R_a denotes a hydrogen atom,

R_b denotes a 1-phenylethyl group or a phenyl group wherein the phenyl nucleus is substituted by the radicals R_1 and R_2 , whilst

R_1 and R_2 , which may be identical or different, each denote a hydrogen, fluorine, chlorine or bromine atom,

X denotes a nitrogen atom,

A denotes an imino group,

B denotes a carbonyl group,

C denotes a 1,2-vinylene -CH=CH-, ethynylene -C≡C- or 1,3-butadien-1,4-ylene -CH=CH-
CH=CH- group, |

D denotes a methylene group,

E denotes a dimethylamino, diethylamino, Bis(2-methoxyethyl)amino, *N*-methyl-*N*-(2-methoxyethyl)amino, *N*-ethyl-*N*-(2-methoxyethyl)amino, *N*-methyl-*N*-cyclopropylamino, *N*-methyl-*N*-cyclopropylmethyl-amino, *N*-methyl-*N*-(1-methoxy-2-propyl)amino, *N*-methyl-*N*-(2-methoxypropyl)amino or *N*-methyl-*N*-(3-methoxypropyl)amino, *N*-methyl-*N*-(tetrahydrofuran-3-yl)amino, *N*-methyl-*N*-(tetrahydropyran-4-yl)amino group,

R_c denotes a cyclopropylmethoxy, cyclobutyloxy or cyclopentyloxy group,

or the tautomers, or stereoisomers or pharmaceutically acceptable salts thereof.

6. (previously presented) The following compounds of general formula I according to claim 1:

4-[(3-Chloro-4-fluorophenyl)amino]-6-{[4-(*N,N*-diethylamino)-1-oxo-2-buten-1-yl]amino}-7-cyclopropylmethoxyquinazoline

or pharmaceutically acceptable salts thereof.

7. (previously presented) Pharmaceutically acceptable salts of the compounds according to one of claims 1 to 6 with inorganic or organic acids or bases.

8. (currently amended) Pharmaceutical composition containing a compound according to at least one of claims 1 to 6, optionally together with ~~one or more inert carriers and/or diluents~~ an inert carrier and with or without a diluent.

9. (currently amended) A method for treating ~~or preventing~~ a disease comprising administering a pharmaceutical composition according to one of claims 1 to 6, wherein said disease is selected from the group consisting of: ~~benign or~~ malignant tumors, diseases of the airways and lungs and diseases of the gastrointestinal tract and the bile duct and gall bladder.

Claims 10-11 (canceled)

12. (currently amended) Pharmaceutical compositions containing a physiologically acceptable salt according to claim 7, optionally together with ~~one or more inert carriers and/or diluents~~ an inert carrier and with or without a diluent.

13. (currently amended) A method for treating ~~or preventing~~ a disease comprising administering a pharmaceutical composition according to claim 7, wherein said disease is selected from the group consisting of: ~~benign or malignant tumors, diseases of the airways and lungs and diseases of the gastrointestinal tract and the bile duct and gall bladder.~~